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INDEX OF SUBJECTS.

TRANSACTIONS. 1918.

Single organic compounds of known empirical formula will be found in the Formula Index, p. 1004.

FORMULA INDEX.

THE following index of organic compounds of known empirical formula is arranged according to Richter's system (see *Lexikon der Kohlenstoff-Verbindungen*).

The elements are given in the order C, H, O, N, Cl, Br, I, F, S, P, and the remainder alphabetically.

The compounds are arranged—

Firstly, in groups according to the number of carbon atoms (thus C_1 group, C_2 group, etc.).

Secondly, according to the number of other elements besides carbon contained in the molecule (thus 5 IV indicates that the molecule contains five carbon atoms and four other elements).

Thirdly, according to the nature of the elements present in the molecule (given in the above order).

Fourthly, according to the number of atoms of each single element (except carbon) present in the molecule.

Salts are placed with the compounds from which they are derived. The chlorides, bromides, iodides, and cyanides of quaternary ammonium bases, however, are registered as group-substances.

C_1 Group.

C_2 Group.

C_3 Group.

C_4 Group.

4 III

C_5 Group.

C_6 Group.

6 II

6 III

C_7 Group.

7 III

7 IV

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C_8 Group.

$C_8H_6O_4$ Terephthalic acid, menthyl alkyl esters of (COHEN and DE PENNINGTON), 57.

$C_8H_8O_2$ Phenylacetic acid, association of, in benzene (INNES), 432.

$C_8H_8O_3$ Methyl salicylate, association of, in benzene (INNES), 431.

$C_8H_{11}N$ *p*-Methylbenzylamine, stannochloride of (DRUCE), 718.

8 III

$C_8H_5O_6N$ Nitroterephthalic acid, menthyl alkyl esters of (COHEN and DE PENNINGTON), 57.

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$C_8H_7O_8N_3$ 3:5:6-Trinitroveratrole (POLLECOFF and ROBINSON), 654.

$C_8H_8O_3N_2$ *p*-Nitromethylformanilide (MORGAN and GRIST), 690.

$C_8H_8O_6N_2$ 5:6-Dinitroguaiacol methyl ether (POLLECOFF and ROBINSON), 651.

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Acetophenoneoxime, association of, in benzene (INNES), 432.

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$C_8H_9O_2Br$ 3-Bromoveratrole (SIMONSEN and RAU), 785.

$C_8H_9O_3N$ 6-Nitro-*o*-tolyl methyl ether (SIMONSEN), 781.

$C_8H_{10}ON_2$ *p*-Aminomethylformanilide (MORGAN and GRIST), 691.

$C_8H_{10}O_4N_2$ 5-Nitro-4-aminoveratrole (SIMONSEN and RAU), 27.

8 IV

$C_8H_5ON_3Cl_2$ 4-Acetyldichloro-*p*-phenylene-1-diazo-4-imides (MORGAN and CLEAGE), 594.

$C_8H_8O_2NCl$ *p*-Chloromandelamide, and its salts (RULE), 17.

$C_8H_{10}O_2NBr$ Bromoaminoveratroles (SIMONSEN and RAU), 786.

C_9 Group.

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$C_9H_{10}O_3$ Ethyl *p*-hydroxybenzoate, association of, in benzene (INNES), 434.

2-Methoxy-*m*-toluic acid, and its silver salt (SIMONSEN), 779.

9 III

$C_9H_9O_2Br$ α-Bromo-β-phenylpropionic acid, sign of the product in conversion of, into α-amino-β-phenylpropionic acid (SENTER, DREW, and MARTIN), 151.

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9 IV

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10 II

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$C_{10}H_8S_2$ β-Naphthylthiosulphoxylic acid, and its sodium salt (WHITE), 608.

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10 III

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$C_{10}H_{12}O_6N_2$ 5-Nitro-4-acetylaminoveratrole (SIMONSEN and RAU), 27.

$C_{10}H_{13}O_2N_3$ 2-Methoxy-*m*-tolualdehyde semicarbazone (SIMONSEN), 777.

$C_{10}H_{17}ON$ Camphoroxime, association of, in benzene (INNES), 432.

10 IV

$C_{10}H_9O_4N_2Cl$ 1-Chloro-2:4-dinitro-*ar*-tetrahydronaphthalene (GREEN and ROWE), 972.

C_{11} Group.

11 III

11 IV

C_{12} Group.

12 III

C_{13} Group.

$C_{13}H_{10}O$ Benzophenone, association of, in benzene (INNES), 430.

$C_{13}H_{12}N_2$ *p*-Aminoacetophenonepiperylhydrazone(+H_2O) (WEINHAGEN), 587.

$C_{13}H_{24}O_{11}$ Methyldigalactoside (CUNNINGHAM), 602.

13 III

$C_{13}H_{11}ON$ Formyldiphenylamine, association of, in benzene (INNES), 431.

$C_{13}H_{11}O_2N$ 2:4-Dihydroxybenzylideneaniline (SENIER and GALLAGHER), 33.

$C_{13}H_{14}O_2N_2$ Piperonalpiperylhydrazone (WEINHAGEN), 585.

$C_{13}H_{11}O_5N_3$ 3:5-Dinitroaceto-*n*-butyl-*p*-toluidide (REILLY and HICKINBOTTOM), 990.

$C_{13}H_{18}ON_2$ Anisaldehydepiperylhydrazone (WEINHAGEN), 586.

$C_{13}H_{12}O_5N_2$ Nitroaceto-*n*-butyl-*p*-toluidides (REILLY and HICKINBOTTOM), 989.

$C_{13}H_{17}ON$ Aceto-*n*-butyl-*o*- and -*p*-toluidides (REILLY and HICKINBOTTOM), 979.

13 IV

$C_{13}H_{10}O_2NBr$ 2:4-Dihydroxybenzylidenebromoanilines (SENIER and GALLAGHER), 34.

C_{14} Group.

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14 II

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15 III

$C_{15}H_{10}O_{13}N_4$ 3:5-Dinitro-2-methoxyphenyl carbonate (POLLECOFF and ROBINSON), 649.

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$C_{15}H_{12}NBr$ Cinnamylidenebromoanilines (SENIER and GALLAGHER), 30.

$C_{15}H_{24}O_4N_2$ 2-Nitrodi-*n*-butyl-*p*-toluidine, and its hydrochloride (REILLY and HICKINBOTTOM), 994.

15 IV

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C_{16} Group.

16 III

16 IV

C_{17} Group.

17 III

17 IV

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C_{19} Group.

$C_{19}H_{16}O_4$ 7-Acetoxy-3-benzyl-2-methyl-1:4-benzopyrone (CRABTREE and ROBINSON), 867.

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19 III

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C_{20} Group.

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$C_{20}H_{20}O_4$ Berberidene (PERKIN), 761.

$C_{20}H_{20}O_5$ Ketodihydroberberidene (PERKIN), 763.

$C_{20}H_{28}O_4$ Ethyl menthyl terephthalate (COHEN and DE PENNINGTON), 63.

$C_{20}H_{30}O_2$ Abietic acid, cerous salt (MORRELL), 120.

$C_{20}H_{38}O_{11}$ Heptamethyl methyllactoside (HAWORTH and LEITCH), 195.
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C_{26} Group.

C_{27} Group.

C_{28} Group.

C_{29} Group.

C_{30} Group.

C_{32} Group.

ERRATA.

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243	15*	*for*	"water" *read* "oxygen."
774	15*	,,	"sodium hypochlorite" *read* "potassium permanganate."

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31	19	,,	"$C_{15}H_{15}O_3N_2$" *read* "$C_{16}H_{14}O_2N_2$."
100	3*	,,	"4-*di*-n-*butylaminoazobenzene*-4′-*sulphonic acid*" read "4-*sulphobenzeneazodi*-n-*butylaniline*."
100	1	,,	"4-n-*butylaminoazobenzene*-4′-*sulphonic acid*" read "4-p-*sulphobenzeneazo*-n-*butylaniline*."
122	16*	,,	"$Ce_2(C_{18}H_{31}O_2,O)_3$" *read* "$Ce(C_{18}H_{31}O_2,O_3)_3$."
123	4	,,	"(Average temperature)" *read* "(Room temperature)."
540	4	,,	"+ 50·6°" *read* "— 50·6°."
774	6	,,	"$C_{20}H_{18}O_7$" *read* "$C_{20}H_{14}O_7$."

* From bottom.